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* * * * * * * * * * Welcome to STN International * * * * * * * * * *

NEWS 1		Web Page URLs for STN Seminar Schedule - N. America
NEWS 2		"Ask CAS" for self-help around the clock
NEWS 3	AUG 09	INSPEC enhanced with 1898-1968 archive
NEWS 4	AUG 28	ADISCTI Reloaded and Enhanced
NEWS 5	AUG 30	CA(SM)/CAplus(SM) Austrian patent law changes
NEWS 6	SEP 21	CA/CAplus fields enhanced with simultaneous left and right truncation
NEWS 7	SEP 25	CA(SM)/CAplus(SM) display of CA Lexicon enhanced
NEWS 8	SEP 25	CAS REGISTRY(SM) no longer includes Concord 3D coordinates
NEWS 9	SEP 25	CAS REGISTRY(SM) updated with amino acid codes for pyrrolysine
NEWS 10	SEP 28	CEABA-VTB classification code fields reloaded with new classification scheme
NEWS 11	OCT 19	LOGOFF HOLD duration extended to 120 minutes
NEWS 12	OCT 19	E-mail format enhanced
NEWS 13	OCT 23	Option to turn off MARPAT highlighting enhancements available
NEWS 14	OCT 23	CAS Registry Number crossover limit increased to 300,000 in multiple databases
NEWS 15	OCT 23	The Derwent World Patents Index suite of databases on STN has been enhanced and reloaded
NEWS 16	OCT 30	CHEMLIST enhanced with new search and display field
NEWS 17	NOV 03	JAPIO enhanced with IPC 8 features and functionality
NEWS 18	NOV 10	CA/CAplus F-Term thesaurus enhanced
NEWS 19	NOV 10	STN Express with Discover! free maintenance release Version 8.01c now available
NEWS 20	NOV 20	CAS Registry Number crossover limit increased to 300,000 in additional databases
NEWS 21	NOV 20	CA/CAplus to MARPAT accession number crossover limit increased to 50,000
NEWS 22	DEC 01	CAS REGISTRY updated with new ambiguity codes
NEWS 23	DEC 11	CAS REGISTRY chemical nomenclature enhanced
NEWS 24	DEC 14	WPIDS/WPINDEX/WPIX manual codes updated
NEWS 25	DEC 14	GBFULL and FRFULL enhanced with IPC 8 features and functionality
NEWS 26	DEC 18	CA/CAplus pre-1967 chemical substance index entries enhanced with preparation role
NEWS 27	DEC 18	CA/CAplus patent kind codes updated
NEWS 28	DEC 18	MARPAT to CA/CAplus accession number crossover limit increased to 50,000
NEWS 29	DEC 18	MEDLINE updated in preparation for 2007 reload
NEWS 30	DEC 27	CA/CAplus enhanced with more pre-1907 records

NEWS EXPRESS NOVEMBER 10 CURRENT WINDOWS VERSION IS V8.01c, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 25 SEPTEMBER 2006.

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FILE 'HOME' ENTERED AT 18:39:48 ON 29 DEC 2006

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STRUCTURE FILE UPDATES: 28 DEC 2006 HIGHEST RN 916479-39-5
DICTIONARY FILE UPDATES: 28 DEC 2006 HIGHEST RN 916479-39-5

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TSCA INFORMATION NOW CURRENT THROUGH June 30, 2006

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

=> E
"N-(4-(3-(4-AMINOPHENYL)-4-ISOAZOLY)PHENYL)-N'-(1-(4-FLUOROPHENYL)ETHYL)THIOUREA"/CN
25
E1 1
N-(4-(3-(4-(MORPHOLIN-4-YL)QUINOLIN-3-YL)ACRYLOYL)PHENYL)OXALAMIC ACID ETHYL ESTER/CN
E2 1
N-(4-(3-(4-ACETYLPIPERAZIN-1-YL)PROPOXY)NAPHTHALEN-1-YL)-3-FLUORO-5-(PIPERIDIN-1-YL) BENZAMIDE/CN
E3 0 -->
N-(4-(3-(4-AMINOPHENYL)-4-ISOAZOLY)PHENYL)-N'-(1-(4-FLUOROPHENYL)ETHYL)THIOUREA/CN
E4 1 N-(4-(3-(4-AMINOPHENYL)PROPYL)-1,3-THIAZOL-2-YL)ACETAMIDE/CN
E5 1
N-(4-(3-(4-BENZYLPIPERIDIN-1-YL)PROPYL)PHENYL)-N-(4-HYDROXYBENZYL)BENZENESULFONAMIDE /CN
E6 1
N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)-3-METHYLNAPHTHALEN-1-YL)ACETAMIDE/CN

```
E7            1
N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)-3-METHYLNAPHTHALEN-1-YL
)ACETAMIDE ACETATE/CN
E8            1
N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)PHENYL)ACETAMIDE/CN
E9            1
N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)PHENYL)ACETAMIDE
ACETATE/CN
E10           1
N-(4-(3-(4-CHLOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXAMIC
ACID/CN
E11           1
N-(4-(3-(4-CHLOROBENZYLCARBAMOYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXAMIC ACID
ETHYL ESTER/CN
E12           1
N-(4-(3-(4-CHLOROPHENOXY)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXAMIC ACID/CN
E13           1
N-(4-(3-(4-CHLOROPHENYL)-1-OXO-2-PROPENYL)PHENYL)-2-METHYL-5-NITROBENZENESULFONAMIDE
/CN
E14           1
N-(4-(3-(4-CHLOROPHENYL)BENZO(C)ISOXAZOL-5-YL)PYRIMIDIN-2-YL)ACETAMIDE/CN
E15           1
N-(4-(3-(4-CHLOROPHENYL)ISOXAZOL-5-YL)THIAZOL-2-YL)-N-(3-(MORPHOLIN-4-YL)PROPYL)THIO
PHENE-2-CARBOXAMIDE/CN
E16           1
N-(4-(3-(4-CYANO-3-TRIFLUOROMETHYLPHENYL)-5,5-DIMETHYL-4-OXO-2-THIOXOIMIDAZOLIDIN-1-
YL)PHENYL)ACETAMIDE/CN
E17           1
N-(4-(3-(4-CYANO-3-TRIFLUOROMETHYLPHENYL)-5,5-DIMETHYL-4-OXO-2-THIOXOIMIDAZOLIDIN-1-
YL)PHENYL)METHANESULFONAMIDE/CN
E18           1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)-2-MORPHOLIN-
4-YL-2-OXOACETAMIDE/CN
E19           1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)-N'-(2-METHOX
YETHYL)OXALAMIDE/CN
E20           1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)-N'-ISOPROPYL
OXALAMIDE/CN
E21           1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)-N'-PROPYLOXA
LAMIDE/CN
E22           1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)MALONAMIC
ACID/CN
E23           1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXALAMIDE/CN
E24           1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXAMIC
ACID/CN
E25           1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXAMIC ACID
ETHYL ESTER/CN

=>
=> E
"N-(4-(3-(4-AMINOPHENYL)-4-ISOxAZOLYl)PHENYL)-N'-(1-(4-FLUOROPHENYL)ETHYL)THIOUREA"/
CN 25
E1            1      N-(4-(3-(4-(MORPHOLIN-4-YL)QUINOLIN-3-YL)ACRYLOYL)PHENYL)OXA
                     LAMIC ACID ETHYL ESTER/CN
E2            1      N-(4-(3-(4-ACETYLPIPERAZIN-1-YL)PROPOXY)NAPHTHALEN-1-YL)-3-F
                     LUORO-5-(PIPERIDIN-1-YL)BENZAMIDE/CN
E3            0 --> N-(4-(3-(4-AMINOPHENYL)-4-ISOXAZOLYL)PHENYL)-N'-(1-(4-FLUORO
                     PHENYL)ETHYL)THIOUREA/CN
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E4              1      N-(4-(3-(4-AMINOPHENYL)PROPYL)-1,3-THIAZOL-2-YL)ACETAMIDE/CN
E5              1      N-(4-(3-(4-BENZYLPIPERIDIN-1-YL)PROPYL)PHENYL)-N-(4-HYDROXYB
                       ENZYL)BENZENESULFONAMIDE/CN
E6              1      N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)
                       -3-METHYLNAPHTHALEN-1-YL)ACETAMIDE/CN
E7              1      N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)
                       -3-METHYLNAPHTHALEN-1-YL)ACETAMIDE ACETATE/CN
E8              1      N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)
                       PHENYL)ACETAMIDE/CN
E9              1      N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)
                       PHENYL)ACETAMIDE ACETATE/CN
E10             1      N-(4-(3-(4-CHLOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)OXAMIC ACID/CN
E11             1      N-(4-(3-(4-CHLOROBENZYLCARBAMOYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)OXAMIC ACID ETHYL ESTER/CN
E12             1      N-(4-(3-(4-CHLOROPHENOXY)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHEN
                       YL)OXAMIC ACID/CN
E13             1      N-(4-(3-(4-CHLOROPHENYL)-1-OXO-2-PROPENYL)PHENYL)-2-METHYL-5
                       -NITROBENZENESULFONAMIDE/CN
E14             1      N-(4-(3-(4-CHLOROPHENYL)BENZO(C)ISOXAZOL-5-YL)PYRIMIDIN-2-YL
                       )ACETAMIDE/CN
E15             1      N-(4-(3-(4-CHLOROPHENYL)ISOXAZOL-5-YL)THIAZOL-2-YL)-N-(3-(MO
                       RPHOLIN-4-YL)PROPYL)THIOPHENE-2-CARBOXAMIDE/CN
E16             1      N-(4-(3-(4-CYANO-3-TRIFLUOROMETHYLPHENYL)-5,5-DIMETHYL-4-OXO
                       -2-THIOXOIMIDAZOLIDIN-1-YL)PHENYL)ACETAMIDE/CN
E17             1      N-(4-(3-(4-CYANO-3-TRIFLUOROMETHYLPHENYL)-5,5-DIMETHYL-4-OXO
                       -2-THIOXOIMIDAZOLIDIN-1-YL)PHENYL)METHANESULFONAMIDE/CN
E18             1      N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)-2-MORPHOLIN-4-YL-2-OXOACETAMIDE/CN
E19             1      N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)-N'-(2-METHOXYETHYL)OXALAMIDE/CN
E20             1      N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)-N'-ISOPROPYLOXALAMIDE/CN
E21             1      N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)-N'-PROPYLOXALAMIDE/CN
E22             1      N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)MALONAMIC ACID/CN
E23             1      N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)OXALAMIDE/CN
E24             1      N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)OXAMIC ACID/CN
E25             1      N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)OXAMIC ACID ETHYL ESTER/CN

=> E
"N1-(4-(3-(4-AMINOPHENYL)-4-ISOxAZOLYl)PHENYL)-N2-(1-(4-FLUOROPHENYL)ETHYL)THIOUREA"
/CN 25
E1              1      N1-(4-(2,4-DIMETHYLTHIAZOL-5-YL)PYRIMIDIN-2-YL)-4-METHOXY-N3
                       ,N3-DIMETHYLBENZENE-1,3-DIAMINE/CN
E2              1      N1-(4-(2-(4-FLUOROPHENYL)-6-TRIFLUOROMETHYLPYRAZOLO(1,5-A)PY
                       RIDIN-3-YL)PYRIMIDIN-2-YL)-N3,N3-DIMETHYLPROPANE-1,3-DIAMINE
                       /CN
E3              0 --> N1-(4-(3-(4-AMINOPHENYL)-4-ISOXAZOLYL)PHENYL)-N2-(1-(4-FLUOR
                       OPHENYL)ETHYL)THIOUREA/CN
E4              1      N1-(4-(3-BROMO-4-((2,4-DIFLUOROBENZYL)OXY)-6-METHYL-2-OXO-2H
                       -PYRIDIN-1-YL)BENZYL)-L-SERINAMIDE HYDROCHLORIDE/CN
E5              1      N1-(4-(3-METHYLCARBAMOYL-4-METHYL-7,8-METHYLENEDIOXY-3,4-DIH
                       YDRO-5H-2,3-BENZODIAZEPIN-1-YL)PHENYL)-N3-METHYLUREA/CN
E6              1      N1-(4-(4-AMINO-7-(4-OXOCYCLOHEXYL)-7H-PYRROLO(2,3-D)PYRIMIDI
                       N-5-YL)-2-FLUOROPHENYL)-2,3-DICHLORO-1-BENZENESULFONAMIDE/CN
E7              1      N1-(4-(4-AMINO-7-(8-METHYL-8-AZA(3.2.1)BICYCLOOCTAN-3-YL)-7H
                       -PYRROLO(2,3-D)PYRIMIDIN-5-YL)-2-FLUOROPHENYL)-2,3-DICHLORO-
                       1-BENZENESULFONAMIDE/CN
E8              1      N1-(4-(4-CHLOROPHENOXY)PHENYL)-N,N-DIMETHYLUREA/CN
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E9              1      N1-(4-(DIHEXYLAMINO)PHENYL)-N1,N4,N4-TRIHEXYL-1,4-PHENYLENED
                       IAMINE/CN
E10             1      N1-(4-AMINOBUTYL)-N4,N4-BIS(2-HYDROXYETHYL)-2-NITRO-P-PHENYL
                       ENEDIAMINE/CN
E11             1      N1-(4-BROMO-2-((2-FLUOROPHENYL)CARBONYL)PHENYL)-L-ALANINAMID
                       E/CN
E12             1      N1-(4-BROMO-2-((2-FLUOROPHENYL)CARBONYL)PHENYL)GLYCINAMIDE/C
                       N
E13             1      N1-(4-BROMOBENZYL)-N1-(PYRID-2-YL)BUTANE-1,4-DIAMINE/CN
E14             1      N1-(4-BROMOPHENYL)-N2-HYDROXY-2-OXO-2-PHENYLACETAMIDINE/CN
E15             1      N1-(4-BUTOXYBENZYL)-5,6-DIHYDROTHIOURACIL/CN
E16             1      N1-(4-CHLOROBENZYL)-4-(3-FLUORO-1-PIPERIDINYL)-4-OXO-1,3-(S)
                       -BUTANEDIAMINE BIS(TRIFLUOROACETATE)/CN
E17             1      N1-(4-CHLOROPHENYL)-N3-CYANOGUANIDINE/CN
E18             1      N1-(4-IMINO-1,3-DIMETHYL-2,6-DIOXOHEXAHYDROPYRIMIDIN-5-YL)SU
                       LFANILAMIDE/CN
E19             1      N1-(4-ISOPROPENYLPHENYL)-N3,N3-DIMETHYLSEMICARBAZIDE/CN
E20             1      N1-(4-ISOPROPOXYBENZOYL)-P-AMINOBENZENESULFONAMIDE/CN
E21             1      N1-(4-ISOPROPOXYBENZOYL)SULFANILAMIDE/CN
E22             1      N1-(4-METHYL-2-PYRIDYL)SULFANILAMIDE SODIUM/CN
E23             1      N1-(4-METHYL-2-PYRIDYL-6-TRIFLUOROMETHYL)SULFANILAMIDE/CN
E24             1      N1-(4-METHYL-2-PYRIMIDINYL)SULFANILAMIDE/CN
E25             1      N1-(4-METHYL-2-THIAZOLYL)SULFANILAMIDE/CN

=> E
"N-(4-(3-(4-AMINOPHENYL)-4-ISOxAZOLYl)PHENYL)-N-(1-(4-FLUOROPHENYL)ETHYL)THIOUREA"/C
N 25
E1              1      N-(4-(3-(4-(MORPHOLIN-4-YL)QUINOLIN-3-YL)ACRYLOYL)PHENYL)OXA
                       LAMIC ACID ETHYL ESTER/CN
E2              1      N-(4-(3-(4-ACETYLPIPERAZIN-1-YL)PROPOXY)NAPHTHALEN-1-YL)-3-F
                       LUORO-5-(PIPERIDIN-1-YL)BENZAMIDE/CN
E3              0 --> N-(4-(3-(4-AMINOPHENYL)-4-ISOXAZOLYL)PHENYL)-N-(1-(4-FLUOROP
                       HENYL)ETHYL)THIOUREA/CN
E4              1      N-(4-(3-(4-AMINOPHENYL)PROPYL)-1,3-THIAZOL-2-YL)ACETAMIDE/CN
E5              1      N-(4-(3-(4-BENZYLPIPERIDIN-1-YL)PROPYL)PHENYL)-N-(4-HYDROXYB
                       ENZYL)BENZENESULFONAMIDE/CN
E6              1      N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)
                       -3-METHYLNAPHTHALEN-1-YL)ACETAMIDE/CN
E7              1      N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)
                       -3-METHYLNAPHTHALEN-1-YL)ACETAMIDE ACETATE/CN
E8              1      N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)
                       PHENYL)ACETAMIDE/CN
E9              1      N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)
                       PHENYL)ACETAMIDE ACETATE/CN
E10             1      N-(4-(3-(4-CHLOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)OXAMIC ACID/CN
E11             1      N-(4-(3-(4-CHLOROBENZYLCARBAMOYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)OXAMIC ACID ETHYL ESTER/CN
E12             1      N-(4-(3-(4-CHLOROPHENOXY)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHEN
                       YL)OXAMIC ACID/CN
E13             1      N-(4-(3-(4-CHLOROPHENYL)-1-OXO-2-PROPENYL)PHENYL)-2-METHYL-5
                       -NITROBENZENESULFONAMIDE/CN
E14             1      N-(4-(3-(4-CHLOROPHENYL)BENZO(C)ISOXAZOL-5-YL)PYRIMIDIN-2-YL
                       )ACETAMIDE/CN
E15             1      N-(4-(3-(4-CHLOROPHENYL)ISOXAZOL-5-YL)THIAZOL-2-YL)-N-(3-(MO
                       RPHOLIN-4-YL)PROPYL)THIOPHENE-2-CARBOXAMIDE/CN
E16             1      N-(4-(3-(4-CYANO-3-TRIFLUOROMETHYLPHENYL)-5,5-DIMETHYL-4-OXO
                       -2-THIOXOIMIDAZOLIDIN-1-YL)PHENYL)ACETAMIDE/CN
E17             1      N-(4-(3-(4-CYANO-3-TRIFLUOROMETHYLPHENYL)-5,5-DIMETHYL-4-OXO
                       -2-THIOXOIMIDAZOLIDIN-1-YL)PHENYL)METHANESULFONAMIDE/CN
E18             1      N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)-2-MORPHOLIN-4-YL-2-OXOACETAMIDE/CN
E19             1      N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                       THYLPHENYL)-N'-(2-METHOXYETHYL)OXALAMIDE/CN
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E20          1     N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                   THYLPHENYL)-N'-ISOPROPYLOXALAMIDE/CN
E21          1     N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                   THYLPHENYL)-N'-PROPYLOXALAMIDE/CN
E22          1     N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                   THYLPHENYL)MALONAMIC ACID/CN
E23          1     N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                   THYLPHENYL)OXALAMIDE/CN
E24          1     N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                   THYLPHENYL)OXAMIC ACID/CN
E25          1     N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIME
                   THYLPHENYL)OXAMIC ACID ETHYL ESTER/CN

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L1      STRUCTURE UPLOADED

=> s l1
SAMPLE SEARCH INITIATED 19:25:26 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -        0 TO ITERATE

100.0% PROCESSED        0 ITERATIONS                                0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:            0 TO        0
PROJECTED ANSWERS:               0 TO        0

L2            0 SEA SSS SAM L1
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Connecting via Winsock to STN

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NEWS 4	AUG 28	ADISCTI Reloaded and Enhanced
NEWS 5	AUG 30	CA(SM)/CAplus(SM) Austrian patent law changes
NEWS 6	SEP 21	CA/CAplus fields enhanced with simultaneous left and right truncation
NEWS 7	SEP 25	CA(SM)/CAplus(SM) display of CA Lexicon enhanced
NEWS 8	SEP 25	CAS REGISTRY(SM) no longer includes Concord 3D coordinates
NEWS 9	SEP 25	CAS REGISTRY(SM) updated with amino acid codes for pyrrolysine
NEWS 10	SEP 28	CEABA-VTB classification code fields reloaded with new classification scheme
NEWS 11	OCT 19	LOGOFF HOLD duration extended to 120 minutes
NEWS 12	OCT 19	E-mail format enhanced
NEWS 13	OCT 23	Option to turn off MARPAT highlighting enhancements available
NEWS 14	OCT 23	CAS Registry Number crossover limit increased to 300,000 in multiple databases
NEWS 15	OCT 23	The Derwent World Patents Index suite of databases on STN has been enhanced and reloaded
NEWS 16	OCT 30	CHEMLIST enhanced with new search and display field
NEWS 17	NOV 03	JAPIO enhanced with IPC 8 features and functionality
NEWS 18	NOV 10	CA/CAplus F-Term thesaurus enhanced
NEWS 19	NOV 10	STN Express with Discover! free maintenance release Version 8.01c now available
NEWS 20	NOV 20	CAS Registry Number crossover limit increased to 300,000 in additional databases
NEWS 21	NOV 20	CA/CAplus to MARPAT accession number crossover limit increased to 50,000
NEWS 22	DEC 01	CAS REGISTRY updated with new ambiguity codes
NEWS 23	DEC 11	CAS REGISTRY chemical nomenclature enhanced
NEWS 24	DEC 14	WPIDS/WPINDEX/WPIX manual codes updated
NEWS 25	DEC 14	GBFULL and FRFULL enhanced with IPC 8 features and functionality
NEWS 26	DEC 18	CA/CAplus pre-1967 chemical substance index entries enhanced with preparation role
NEWS 27	DEC 18	CA/CAplus patent kind codes updated
NEWS 28	DEC 18	MARPAT to CA/CAplus accession number crossover limit increased to 50,000
NEWS 29	DEC 18	MEDLINE updated in preparation for 2007 reload
NEWS 30	DEC 27	CA/CAplus enhanced with more pre-1907 records

NEWS EXPRESS NOVEMBER 10 CURRENT WINDOWS VERSION IS V8.01c, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 25 SEPTEMBER 2006.

NEWS HOURS	STN Operating Hours Plus Help Desk Availability
NEWS LOGIN	Welcome Banner and News Items
NEWS IPC8	For general information regarding STN implementation of IPC 8

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 NEWS X25       X.25 communication option no longer available
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=> file registry
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FULL ESTIMATED COST                                           0.21       0.21

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STRUCTURE FILE UPDATES:   28 DEC 2006  HIGHEST RN 916479-39-5
DICTIONARY FILE UPDATES:  28 DEC 2006  HIGHEST RN 916479-39-5

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on property searching in REGISTRY, refer to:

http://www.cas.org/ONLINE/UG/regprops.html

=> E
"N-(4-(3-(4-AMINOPHENYL)-4-ISOAZOLY)PHENYL)-N'-(1-(4-FLUOROPHENYL)ETHYL)THIOUREA"/CN
 25
E1            1
N-(4-(3-(4-(MORPHOLIN-4-YL)QUINOLIN-3-YL)ACRYLOYL)PHENYL)OXALAMIC ACID ETHYL
ESTER/CN
E2            1
N-(4-(3-(4-ACETYLPIPERAZIN-1-YL)PROPOXY)NAPHTHALEN-1-YL)-3-FLUORO-5-(PIPERIDIN-1-YL)
BENZAMIDE/CN
E3            0 -->
N-(4-(3-(4-AMINOPHENYL)-4-ISOAZOLY)PHENYL)-N'-(1-(4-FLUOROPHENYL)ETHYL)THIOUREA/CN
E4            1      N-(4-(3-(4-AMINOPHENYL)PROPYL)-1,3-THIAZOL-2-YL)ACETAMIDE/CN
E5            1
N-(4-(3-(4-BENZYLPIPERIDIN-1-YL)PROPYL)PHENYL)-N-(4-HYDROXYBENZYL)BENZENESULFONAMIDE
/CN
E6            1
N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)-3-METHYLNAPHTHALEN-1-YL
)ACETAMIDE/CN
```

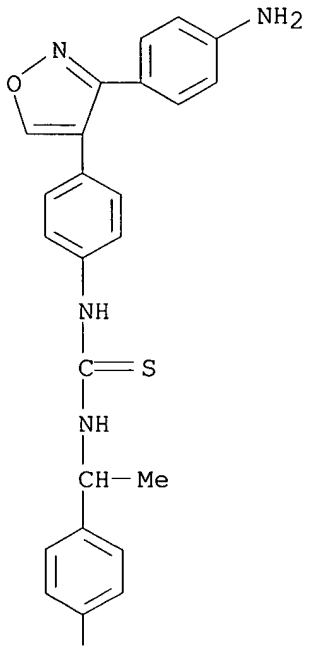
E7 1
N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)-3-METHYLNAPHTHALEN-1-YL)ACETAMIDE ACETATE/CN
E8 1
N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)PHENYL)ACETAMIDE/CN
E9 1
N-(4-(3-(4-CARBAMIMIDOYLPHENYL)-2-OXOOXAZOLIDIN-5-YLMETHOXY)PHENYL)ACETAMIDE ACETATE/CN
E10 1
N-(4-(3-(4-CHLOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXAMIC ACID/CN
E11 1
N-(4-(3-(4-CHLOROBENZYLCARBAMOYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXAMIC ACID ETHYL ESTER/CN
E12 1
N-(4-(3-(4-CHLOROPHENOXY)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXAMIC ACID/CN
E13 1
N-(4-(3-(4-CHLOROPHENYL)-1-OXO-2-PROPENYL)PHENYL)-2-METHYL-5-NITROBENZENESULFONAMIDE/CN
E14 1
N-(4-(3-(4-CHLOROPHENYL)BENZO(C)ISOXAZOL-5-YL)PYRIMIDIN-2-YL)ACETAMIDE/CN
E15 1
N-(4-(3-(4-CHLOROPHENYL)ISOXAZOL-5-YL)THIAZOL-2-YL)-N-(3-(MORPHOLIN-4-YL)PROPYL)THIOPHENE-2-CARBOXAMIDE/CN
E16 1
N-(4-(3-(4-CYANO-3-TRIFLUOROMETHYLPHENYL)-5,5-DIMETHYL-4-OXO-2-THIOXOIMIDAZOLIDIN-1-YL)PHENYL)ACETAMIDE/CN
E17 1
N-(4-(3-(4-CYANO-3-TRIFLUOROMETHYLPHENYL)-5,5-DIMETHYL-4-OXO-2-THIOXOIMIDAZOLIDIN-1-YL)PHENYL)METHANESULFONAMIDE/CN
E18 1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)-2-MORPHOLIN-4-YL-2-OXOACETAMIDE/CN
E19 1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)-N'-(2-METHOXYETHYL)OXALAMIDE/CN
E20 1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)-N'-ISOPROPYLOXALAMIDE/CN
E21 1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)-N'-PROPYLOXALAMIDE/CN
E22 1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)MALONAMIC ACID/CN
E23 1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXALAMIDE/CN
E24 1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXAMIC ACID/CN
E25 1
N-(4-(3-(4-FLUOROBENZENESULFONYL)-4-HYDROXYPHENOXY)-3,5-DIMETHYLPHENYL)OXAMIC ACID ETHYL ESTER/CN

L7 ANSWER 91 OF 177 REGISTRY COPYRIGHT 2007 ACS on STN
RN 736982-21-1 REGISTRY
ED Entered STN: 01 Sep 2004
CN Thiourea, N-[4-[3-(4-aminophenyl)-4-isoxazolyl]phenyl]-N'-[1-(4-fluorophenyl)ethyl]- (9CI) (CA INDEX NAME)
MF C24 H21 F N4 O S
SR CA
LC STN Files: CA, CAPLUS, USPATFULL
DT.CA CAplus document type: Patent
RL.P Roles from patents: BIOL (Biological study); PREP (Preparation); USES (Uses)

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	RID Occurrence Count
C3NO	NOC3	5	C3NO	16.167.5	1
C6	C6	6	C6	46.150.18	3

PAGE 1-A



PAGE 2-A

F

Experimental Property Tags (ETAG)

PROPERTY	NOTE
Mass Spectra	(1) CAS

(1) [redacted] Jonathan David; US 2004157900 A1 2004 CAPLUS

Predicted Properties (PPROP)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	1.0	pH 1 25 deg C	(1)
Bioconc. Factor (BCF)	10.28	pH 2 25 deg C	(1)
Bioconc. Factor (BCF)	53.90	pH 3 25 deg C	(1)
Bioconc. Factor (BCF)	90.90	pH 4 25 deg C	(1)
Bioconc. Factor (BCF)	97.57	pH 5 25 deg C	(1)
Bioconc. Factor (BCF)	98.29	pH 6 25 deg C	(1)
Bioconc. Factor (BCF)	98.36	pH 7 25 deg C	(1)
Bioconc. Factor (BCF)	98.36	pH 8 25 deg C	(1)
Bioconc. Factor (BCF)	98.26	pH 9 25 deg C	(1)
Bioconc. Factor (BCF)	97.29	pH 10 25 deg C	(1)
Boiling Point (BP)	583.4+/-60.0 deg C	760 Torr	(1)
Density (DEN)	1.309+/-0.06 g/cm**3	760 Torr	(1)
Enthalpy of Vap. (HVAP)	87.20+/-3.0 kJ/mol	760 Torr	(1)
Flash Point (FP)	306.6+/-32.9 deg C		(1)
Freely Rotatable Bonds (FRB)	6		(1)
H acceptors (HAC)	5		(1)
H donors (HD)	4		(1)
Hydrogen Donors/Acceptors Sum (HDAS)	9		(1)
Koc (KOC)	8.42	pH 1 25 deg C	(1)
Koc (KOC)	97.08	pH 2 25 deg C	(1)
Koc (KOC)	508.80	pH 3 25 deg C	(1)
Koc (KOC)	858.08	pH 4 25 deg C	(1)
Koc (KOC)	921.04	pH 5 25 deg C	(1)
Koc (KOC)	927.84	pH 6 25 deg C	(1)
Koc (KOC)	928.52	pH 7 25 deg C	(1)
Koc (KOC)	928.49	pH 8 25 deg C	(1)
Koc (KOC)	927.58	pH 9 25 deg C	(1)
Koc (KOC)	918.43	pH 10 25 deg C	(1)
LOGD (LOGD)	0.88	pH 1 25 deg C	(1)
LOGD (LOGD)	1.94	pH 2 25 deg C	(1)
LOGD (LOGD)	2.66	pH 3 25 deg C	(1)
LOGD (LOGD)	2.89	pH 4 25 deg C	(1)
LOGD (LOGD)	2.92	pH 5 25 deg C	(1)
LOGD (LOGD)	2.92	pH 6 25 deg C	(1)
LOGD (LOGD)	2.92	pH 7 25 deg C	(1)
LOGD (LOGD)	2.92	pH 8 25 deg C	(1)
LOGD (LOGD)	2.92	pH 9 25 deg C	(1)
LOGD (LOGD)	2.92	pH 10 25 deg C	(1)
LOGP (LOGP)	2.925+/-0.555	25 deg C	(1)
Mass Intrinsic Solubility (ISLB.MASS)	0.0069 g/L	25 deg C	(1)
Mass Solubility (SLB.MASS)	0.78 g/L	pH 1 25 deg C	(1)
Mass Solubility (SLB.MASS)	0.069 g/L	pH 2 25 deg C	(1)
Mass Solubility (SLB.MASS)	0.013 g/L	pH 3 25 deg C	(1)
Mass Solubility (SLB.MASS)	0.0078 g/L	pH 4 25 deg C	(1)
Mass Solubility (SLB.MASS)	0.0074 g/L	pH 5 25 deg C	(1)
Mass Solubility (SLB.MASS)	0.0069 g/L	pH 6 25 deg C	(1)
Mass Solubility (SLB.MASS)	0.0069 g/L	pH 7 25 deg C	(1)
Mass Solubility (SLB.MASS)	0.0069 g/L	pH 8 25 deg C	(1)
Mass Solubility (SLB.MASS)	0.0069 g/L	pH 9 25 deg C	(1)
Mass Solubility (SLB.MASS)	0.0074 g/L	pH 10 25 deg C	(1)
Mass Solubility (SLB.MASS)	0.0069 g/L	Unbuffered Water	(1)

		pH 7.00 25 deg C	
Molar Intrinsic Solubility (ISLB.MOL)	0.000016 mol/L	25 deg C	(1)
Molar Solubility (SLB.MOL)	0.0018 mol/L	pH 1 25 deg C	(1)
Molar Solubility (SLB.MOL)	0.00016 mol/L	pH 2 25 deg C	(1)
Molar Solubility (SLB.MOL)	0.000030 mol/L	pH 3 25 deg C	(1)
Molar Solubility (SLB.MOL)	0.000018 mol/L	pH 4 25 deg C	(1)
Molar Solubility (SLB.MOL)	0.000017 mol/L	pH 5 25 deg C	(1)
Molar Solubility (SLB.MOL)	0.000016 mol/L	pH 6 25 deg C	(1)
Molar Solubility (SLB.MOL)	0.000016 mol/L	pH 7 25 deg C	(1)
Molar Solubility (SLB.MOL)	0.000016 mol/L	pH 8 25 deg C	(1)
Molar Solubility (SLB.MOL)	0.000016 mol/L	pH 9 25 deg C	(1)
Molar Solubility (SLB.MOL)	0.000017 mol/L	pH 10 25 deg C	(1)
Molar Solubility (SLB.MOL)	0.000016 mol/L	Unbuffered Water pH 7.00 25 deg C	(1)
Molar Volume (MVOL)	330.2+/-3.0 cm**3/mol	20 deg C 760 Torr	(1)
Molecular Weight (MW)	432.51		(1)
PKA (PKA)	12.00+/-0.70	Most Acidic 25 deg C	(1)
PKA (PKA)	2.91+/-0.10	Most Basic 25 deg C	(1)
Polar Surface Area (PSA)	108.20 A**2		(1)
Vapor Pressure (VP)	1.34E-13 Torr	25 deg C	(1)

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software V8.19 ((C) 1994-2007 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1

AN 141:190783 CA
TI Preparation of isoxazole-containing thiourea inhibitors useful for treatment of varicella zoster virus
IN Bloom, Jonathan David
PA Wyeth Holdings Corporation, USA
SO U.S. Pat. Appl. Publ., 9 pp.
CODEN: USXXCO
DT Patent
LA English
IC ICM C07D261-02
ICS A61K031-42
NCL 514378000
CC 28-6 (Heterocyclic Compounds (More Than One Hetero Atom))
Section cross-reference(s): 1, 63
FAN.CNT 1

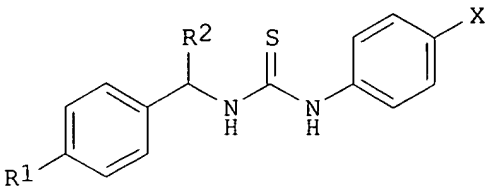
	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 2004157900	A1	20040812	US 2004-772799	20040205
	WO 2004072052	A2	20040826	WO 2004-US3725	20040209
	WO 2004072052	A3	20041111		

W: AE, AE, AG, AL, AL, AM, AM, AM, AT, AT, AU, AZ, AZ, BA, BB, BG, BG, BR, BR, BW, BY, BY, BZ, BZ, CA, CH, CN, CN, CO, CO, CR, CR, CU, CU, CZ, CZ, DE, DE, DK, DK, DM, DZ, EC, EC, EE, EE, EG, ES, ES, FI, FI, GB, GD, GE, GE, GH, GM, HR, HR, HU, HU, ID, IL, IN, IS, JP, JP, KE, KE, KG, KG, KP, KP, KP, KR, KR, KZ, KZ, KZ, LC, LK, LR, LS, LS, LT, LU, LV, MA, MD, MD, MG, MK, MN, MW, MX, MX, MZ, MZ, NA, NI

RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

PRAI US 2003-446602P 20030211

GI



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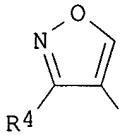

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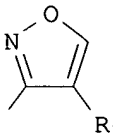

III

AB The title compds. [I; R1 = halo, H; R2 = alkyl; X = II, III; R3 = alkyl, cycloalkyl, hydroxymethyl, etc.; R4 = alkyl which may be further substituted with (un)substituted Ph, cycloalkyl, pyridyl, etc.], useful for inhibiting replication of a herpes virus, were prepared E.g., a multi-step synthesis of 1-[4-(4-benzylisoxazol-3-yl)phenyl]-3-[1-(4-fluorophenyl)ethyl]thiourea (IV), was given. Seventeen title compds. I were prepared as described for IV, and tested for activity as herpes virus inhibitors (IC50 values against VZV, MTS, CMV, HSV and RSV were given).

ST isoxazolylphenyl benzyl thiourea prepn antiviral herpes virus; varicella zoster virus isoxazolylphenyl benzyl thiourea prepn

IT Antiviral agents
Human
Human herpesvirus
Human herpesvirus 3
Human herpesvirus 5
(preparation of isoxazole-containing thiourea inhibitors useful for treatment of
varicella zoster virus)

IT Infection
(viral; preparation of isoxazole-containing thiourea inhibitors useful for treatment of varicella zoster virus)

IT 736982-12-0P 736982-13-1P 736982-14-2P 736982-15-3P 736982-16-4P
736982-17-5P 736982-18-6P 736982-19-7P 736982-20-0P 736982-21-1P
736982-22-2P 736982-23-3P 736982-24-4P 736982-25-5P 736982-26-6P
736982-27-7P 736982-28-8P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of isoxazole-containing thiourea inhibitors useful for treatment of
varicella zoster virus)

IT 100-10-7, 4-(Dimethylamino)benzaldehyde 623-04-1, 4-Aminobenzyl alcohol 10147-11-2, 3-Phenyl-1-propyne 14235-81-5, 4-Ethynylaniline 182565-27-1
RL: RCT (Reactant); RACT (Reactant or reagent)

(preparation of isoxazole-containing thiourea inhibitors useful for treatment of
varicella zoster virus)
IT 2929-84-2P, 4-Dimethylaminobenzaldehyde oxime 144072-29-7P
144072-30-0P 157991-82-7P 170727-03-4P 190446-52-7P 736982-09-5P
736982-10-8P 736982-11-9P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation of isoxazole-containing thiourea inhibitors useful for treatment of
varicella zoster virus)